

Treebeard : An Optimizing Compiler for Tree Based ML Inference

ABSTRACT

Decision tree ensembles are commonly used machine learning models generated by machine learning techniques like gradient boosting and random forests. These models are used in a wide range of applications and are deployed at scale (run repeatedly on billions of data items). Several libraries such as XGBoost, LightGBM, and Sklearn expose algorithms for both training and inference with decision tree ensembles. While these libraries incorporate a limited set of hardware-specific optimizations, they do not specialize the inference code to the model being used, leaving significant performance on the table.

This paper presents a compiler-based approach that automatically generates efficient code for decision tree inference. It develops Treebeard, an extensible compiler, which progressively lowers inference code to LLVM IR through multiple intermediate abstractions. By applying model-specific optimizations at the higher levels, loop optimizations at the middle level, and machine-specific optimizations lower down, Treebeard can specialize inference code for each model on each supported hardware target. To improve model inference performance, Treebeard performs several novel optimizations such as tree tiling, tree walk unrolling, and tree walk interleaving.

We implement Treebeard using the MLIR compiler infrastructure and demonstrate the utility of Treebeard by evaluating it on a diverse set of tree ensemble models. Experimental evaluation demonstrates that Treebeard optimizations improve average latency over a batch of inputs by 2.2X over our baseline version. Further, Treebeard is significantly faster than XGBoost and Treelite in both single-core (2.8X and 5.1X respectively) and multi-core (3.2X and 2.6X respectively) settings.

1. INTRODUCTION

Intro tex here!

2. COMPILER OVERVIEW

Figure 1 shows the high level structure of Treebeard. The input to Treebeard is a serialized decision tree ensemble. Popular frameworks like XGBoost and LightGBM are supported and the system is easily extensible to other frameworks. Given an input model our compiler generates optimized inference code. Specifically it generates a callable batch inference function `predictForest` that, given an array of input rows, computes an array of model predictions.

Treebeard specializes the code generated for inference by progressively optimizing and lowering a high level representation of the `predictForest` function down to LLVM IR [1]. By *lowering*, we mean the process of transforming an operation at a higher abstraction to a sequence of operations at a lower abstraction. Optimizations in Treebeard are implemented using a combination of annotation and lowering. An operation at a higher abstraction is annotated with attributes that indicate what kind of optimization is to be performed while lowering it. The lowering transformation uses this information to generate optimized lower level code. For example, tree tiling and loop order are decided at the highest abstraction. These decisions are communicated to the lowering pass which explicitly encodes them in the lowered IR as shown in figure 2.

Treebeard’s IR has three abstractions as shown in Figure 2. At the highest level (HIR), the input model is represented as a collection of binary trees. This is shown in the second row of figure 2. At this level of abstraction, Treebeard tiles nodes together to transform a binary tree into an n -ary tree and decides what order trees are to be traversed in. In the example in the figure, trees are tiled with a tile size of 2. Tiling is indicated using colored ellipses drawn around nodes that are in the same tile. Tree reordering is another transformation that is performed at this abstraction. One objective of tree reordering is to group identically structured trees so that they can share the same traversal code. With the tiling shown in Figure 2, Tree1 and Tree3 have depth 2, while Tree2 has a depth of 3. Therefore, the trees are reordered so that Tree1 and Tree3 are together.

The high level IR is then lowered to a mid-level IR (MIR). The aim of MIR is to allow optimizations that are independent of the final memory layout of the model. The lowering to MIR performs loop transformations like loop tiling, permutation, parallelization and unrolling. At this level of the IR (the third row in Figure 2), the order in which trees, input row pairs are walked is explicitly represented in the IR using loop nests. Figure 2 shows two possible ways that loop nests could be generated – the first walks all rows for one tree before moving to the next tree while the second walks all trees for one row before moving to the next row. Another aspect handled by this lowering is the fissioning of loops to make sure that trees with the same structure share code. Here, both versions of MIR shown ensure that Tree1 and Tree3 share the same traversal code, while traversal code for Tree2 is different. Additionally, tree walk optimizations, such as tree walk unrolling (shown in figure 2) and peeling are performed on the MIR.

Treebeard then further lowers the IR to explicitly represent

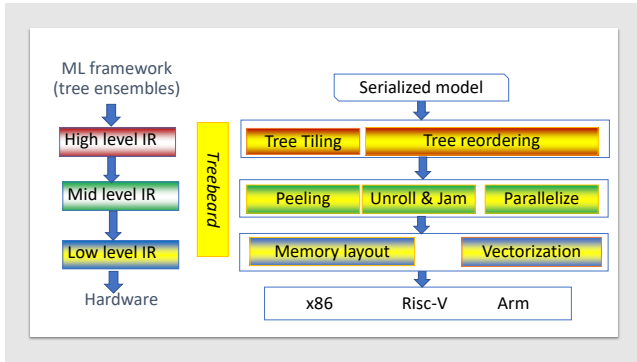


Figure 1: Treebeard compiler structure

the memory layout of the model. Buffers to hold model values are inserted into the generated code and all tree operations in the mid-level IR are lowered to explicitly reference these buffers. Additionally, at this level of the IR, Treebeard generates vectorized code to take advantage of SIMD instructions where applicable. Finally, the inference function is translated to LLVM IR and JIT compiled to executable code.

3. HIGH LEVEL IR OPTIMIZATIONS

This section describes tiling and tree ordering, two optimizations performed at the highest level of abstraction. Recall that here the `predictForest` operator is abstractly represented by just a set of trees.

3.1 Notation

TODO Notation needs to be introduced in the background section, Tiling should be the first subsection We represent a decision tree by $T = (V, E, r)$ where V is the set of nodes, E the set of edges and $r \in V$ is the root. For each node $n \in V$, we define the following.

1. $threshold(n) \in \mathbb{R}$, the threshold value for n .
2. $featureIndex(n) \in \mathbb{N}$, the feature index for n .
3. $left(n) \in V$, the left child of n or \emptyset if n is a leaf. If $left(n) \neq \emptyset$, then $(n, left(n)) \in E$.
4. $right(n) \in V$, the right child of n or \emptyset if n is a leaf. If $right(n) \neq \emptyset$, then $(n, right(n)) \in E$.

We use $L \subseteq V$ to denote the set of leaves.

3.2 Tiling

While a decision tree is naturally represented by a binary tree, it is not the best representation for tree traversal as it (i) requires many memory accesses, (ii) has poor branching structure and (iii) cannot make use of vector instructions. This section proposes a tiling optimization where we group multiple nodes in a decision tree into a single tile, effectively transforming a binary tree into an n -ary tiled tree. This not only allows the compiler to generate vectorized code (see Section 5.1) to traverse trees but also enables spatial locality improvements by grouping nodes that are likely to be accessed together. We demonstrate this with two tiling heuristics later in the section.

Once trees are tiled Treebeard generates tree walks with the code structure shown below.

```

1 ResultType Prediction_Function(...) {
2   // ...
3   Tile t = getRootTile(tree)
4   while (!isLeaf(tree, t)) do {
5     // Evaluate conditions of all nodes in the tile
6     predicates = evaluateTilePredicates(t, rows[i])
7
8     // Move to the correct child of the current tile
9     t = getChildTile(tree, t, predicates)
10  }
11  treePrediction = getLeafValue(t)
12  // ...
13 }
```

The code is just an abstract representation of a tiled tree walk that enables efficient lowering of specific steps in subsequent stages. `evaluateTilePredicates` (speculatively) computes the predicates on all nodes in a tile (line 6). Then `getChildTile` (line 9), uses the computed predicate values to determine which child of the current tile to move to. We defer a description of how these operators are lowered to a later section but focus on tiling algorithms in this section.

Conditions for Efficient Tiling

While any arbitrary partitioning of the nodes of a tree could be considered for tiling we impose a few intuitive constraints. **TODO** `kr : replace n_t with sz ?` Given a $tree = (V, E, r)$ and a tile size n_t we impose the following constraints on the generated tiles $\{T_1, T_2, \dots, T_m\}$.

Partitioning $T_1 \cup T_2 \dots \cup T_m = V$ and $T_i \cap T_j = \emptyset$ for all $i \neq j$

Connectedness If $u, v \in T_i$, there is a (undirected) path connecting u and v fully contained in T_i .

Leaf separation $\forall l \in L : l \in T_i \rightarrow v \notin T_i \ \forall v \in V \setminus \{l\}$

Maximal tiling if there are tiles such that. $|T_i| < n_t$, then there is no $v \in V \setminus \{T_i \cup L\}$ such that $(u, v) \in E$ for some $u \in T_i$.

The **partitioning** and **maximal tiling** constraints together ensure that we group nodes into as few tiles as possible. **Connectedness** ensures that each tile is a sub-tree, a natural grouping of nodes that are likely to be accessed together. The **Leaf separation** constraint ensures that leafs are not tiled along with internal nodes. Leafs in a decision tree need special handling, they are used to check for walk termination and to determine the output (prediction). This constraint ensures that tiles are homogenous, this in-turn allows us to specialize the in-memory layout of trees and also simplifies code generation. We discuss leaf handling and tree layout in Section ?? . We refer to any tiling that satisfies the above constraints as a *valid* tiling.

3.3 Basic Tiling Algorithm

Algorithm 1 shows a simple tiling algorithm that satisfies all the above constraints. Tiling starts at the root and constructs a tile $Tile$ by performing a level order traversal. The call `LevelOrderTraversal(r, n_t)` picks the next n_t nodes according to the standard level order tree traversal algorithm (not described here) applied only to internal nodes of the tree. Once the current tile is constructed, the tiling procedure is recursively performed on all nodes that are destinations for edges going out of the constructed tile. It is easy to see that

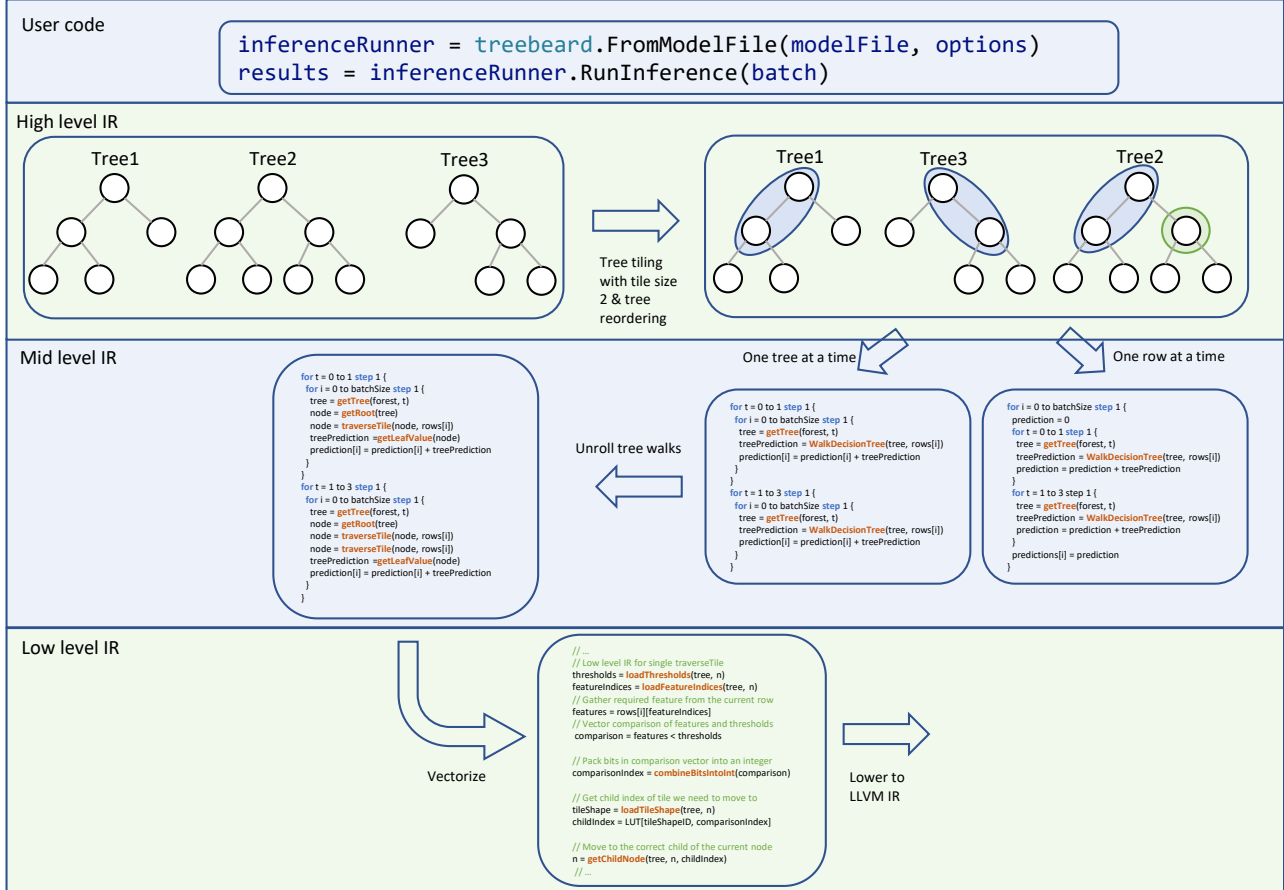


Figure 2: Treebeard IR lowering and optimization details. The three abstraction level in Treebeard’s IR are shown. The high level IR is a tree based IR to perform model level optimization, the mid-level IR is for loop optimizations that are independent of memory layout and the low level IR allows us to perform vectorization and other memory layout dependent optimizations.

the set of tiles constructed by algorithm 1 satisfies the constraints above. **TODO** Pass a modified tree without leafs to the level order traversal call. Explicitly add leaves as separate tiles

Algorithm 1 Uniform tree tiling

```

procedure TILETREE( $tree = (V, E, r), n_t$ )
  if  $r \in L$  then
    return  $\{r\}$ 
  end if
  //Level order traversal to collect  $n_t$  or fewer nodes.
  //Leaves are not included in the constructed tile.
   $Tile \leftarrow \text{LevelOrderTraveral}(r, n_t)$ 
   $Tiles = \{Tile\}$ 
  for  $(u, v) \in \text{Out}(Tile)$  do
     $Tiles \leftarrow Tiles \cup \text{TileTree}(S_v, n_t)$ 
  end for
  return  $Tiles$ 
end procedure

```

One interesting property of this tiling algorithm is that it naturally reduces the imbalance in trees, especially at large tile sizes. As the algorithm traverses down to sparser level of the tree, it naturally groups sub-trees containing chains of nodes, thus balancing the trees. While its possible to further enhance the algorithm to explicitly balance tiled trees, we find that basic tiling suffices in practice.

3.4 Probability Based Tiling

The next algorithm we propose exploits the inherent biases among the leaves of a decision tree. In typical machine learning models some leafs (equivalently outcomes or predictions) are more likely to be reached than others. In such settings, having balanced tiled trees is not sufficient to minimize expected inference time.

Consider for example two machine learning models `airline-ohe` and `epsilon` (also used in our evaluation). Consider the graphs shown in figures 3 and 4 that are generated from training data. Each line in these graphs corresponds to a fixed fraction of the input (say f). A point on a line at coordinate (x, y) means that a fraction y of trees in the model could cover a fraction f of all training inputs with a fraction x of leaves. For example, the first point on the $f = 0.9$ line in figure 3 says that about 52% of trees (y value) need only 1% of their leaves (x value) to cover 90% of the training input. In general, Figure 3 shows that very few leaves are needed to cover a very large fraction of inputs for the benchmark `airline-ohe`. This means that a small fraction of leaves are very likely. We call trees with a small number of extremely likely leaves *leaf biased*.

On the other hand, for the benchmark `epsilon`, figure 4 shows that a trees need a much larger fraction of their leaves to cover a significant fraction of the training input. This means that most trees in the `epsilon` model are not leaf biased.

3.4.1 The Optimization Problem

Observe that the latency of one tree walk is proportional to the number of tiles that need to be evaluated to reach the leaf. It is easy to see that for a leaf biased tree, basic tiling

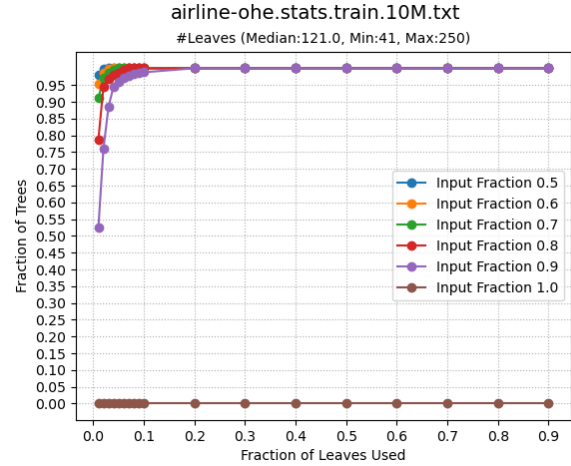


Figure 3: Statistical profile for airline-ohe

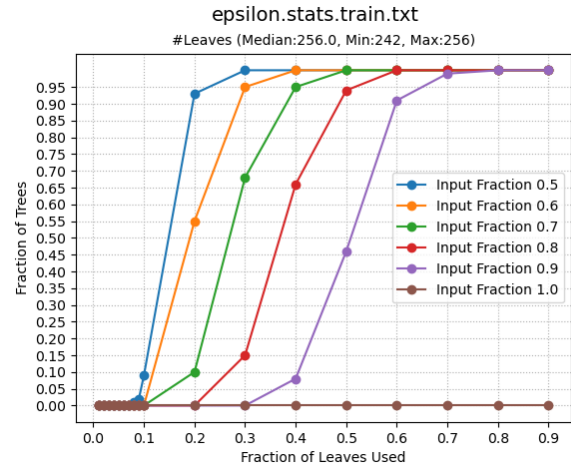


Figure 4: Statistical profile for epsilon

does not optimize for this objective, it considers all leafs to be equally likely.

The goal of probabilistic tiling is to minimize the average inference latency, or equivalently the minimize the expected number of tiles that are evaluated to compute one tree prediction. More formally, the problem is to find a *valid* (as defined in Section 3.2) tiling \mathcal{T} such that the following objective is minimized.

$$\min_{\mathcal{T} \in \mathcal{C}(T)} \sum_{l \in L} p_l \cdot \text{depth}_{\mathcal{T}}(l)$$

where the minimization is over all valid tilings \mathcal{T} of the tree T , $\text{depth}_{\mathcal{T}}(l)$ is the depth of the leaf l given tiling \mathcal{T} . p_l is the probability of reaching leaf l as observed during training.

The above optimization problem can be solved optimally using dynamic programming. We leave this out in the interest of space. Instead, we use the simple greedy algorithm listed in algorithm 2 to construct a valid tiling given the node probabilities¹. The algorithm starts at the root and greedily keeps adding the most probable legal node to the current tile until the maximum tile size is reached. Subsequently, the tiling procedure is recursively performed on all nodes that are destinations for edges going out of the constructed tile.

Algorithm 2 Greedy Probability Based Tree Tiling

```

procedure TILETREE( $T = (V, E, r), n_t$ )
  if  $r \in L$  then
    return  $\{r\}$ 
  end if
   $\text{Tile} \leftarrow \{r\}$ 
  while  $|\text{Tile}| < n_t$  do
     $e = (u, v) \in \text{Out}(\text{Tile})$  st  $p(v)$  is max and  $v \notin L$ 
    if  $e = \emptyset$  then
      break
    end if
     $\text{Tile} = \text{Tile} \cup \{v\}$ 
  end while
   $\text{Tiles} = \{\text{Tile}\}$ 
  for  $(u, v) \in \text{Out}(\text{Tile})$  do
     $\text{Tiles} \leftarrow \text{Tiles} \cup \text{TILETREE}(S_v, n_t)$ 
  end for
  return  $\text{Tiles}$ 
end procedure

```

We find probability based tiling is only beneficial for leaf biased trees². Recall that a tree to be leaf biased if a small fraction of leaves, say α , can cover a large fraction of training inputs, say β . We only perform probability based tiling on trees with thresholds $\alpha = 0.05$ and $\beta = 0.9$ and fall back to uniform tiling otherwise.

3.5 Tree ordering

¹Probabilities for internal nodes can be computed from probabilities for leafs by summing up the probabilities of all leafs that belong to the sub-tree rooted at the internal node. Leaf probabilities are collected during training.

²Turns out that probability based tiling produces many more tile shapes (see Section ??) which directly impacts the cost of `getChildTile` making it more expensive than basic tiling

Specializing the code for each tree in a model comes at a cost. First the code generator needs to generate different code for different trees potentially increasing the size of the generated code. Second some cross tree optimizations (applied at the lower levels of abstraction) like tree walk interleaving require that the multiple trees share identical code. In order to handle this, TREEBEARD pads trees with dummy nodes to make them balanced and then sorts the trees by their depth, so that trees of same depth can share code. Padding is only done for almost balanced trees (as generated by basic tiling), this is ensured by only adding up to a fixed fraction of dummy nodes. **TODO Kr : Can we say something better above. If not can we add a threshold say 10%?. Note : need to mention unrolling of padded sections in MIR.** Once this is done, the loop over the trees is fished so that each of the resulting loops only walks trees of a single depth. Consider for example a forest with 4 trees T_1, T_2, T_3 , and T_4 in that order. Further, assume that T_1 and T_4 have depth 2 while T_2 and T_3 have depth 3. First, Treebeard reorders the trees to be in the order T_1, T_4, T_2, T_3 . Then, the loop over the trees is fished as shown in the following listing.

```

1 forest = ensemble(...)
2 for i = 0 to batchSize step 1 {
3   prediction = 0
4   for t = 0 to 2 step 1 {
5     tree = getTree(forest, t)
6     node = getRoot(tree)
7     node = traverseTreeTile(tree, node, rows[i])
8     treePrediction = getLeafValue(tree, node)
9     prediction = prediction + treePrediction
10  }
11  for t = 2 to 4 step 1 {
12    tree = getTree(forest, t)
13    node = getRoot(tree)
14    node = traverseTreeTile(tree, node, rows[i])
15    node = traverseTreeTile(tree, node, rows[i])
16    treePrediction = getLeafValue(tree, node)
17    prediction = prediction + treePrediction
18  }
19  predictions[i] = prediction
20 }

```

4. LOOP OPTIMIZATIONS AT MID LEVEL

4.1 Tree Walk Interleaving

While tiling and subsequent vectorization gives significant performance gains, profiling the generated code showed that true dependencies between instructions were still causing a significant number of stall cycles. In order to fill these stall cycles, Treebeard does an unroll and jam on the inner most loops of the loop nest. This transformation has the effect of walking multiple tree and input row pairs in an interleaved fashion. The dependency stalls can be hidden by scheduling instructions from independent tree walks in the stall cycles.

This optimization is performed across both the mid-level IR and low-level IR. Loops on which to perform unroll and jam are identified in the mid-level IR. This information is communicated to the lowering passes by annotating these tree walk operations as described in section ???. The lowering passes that transform the mid-level IR to low-level IR interleave operations across independent tree walks.

The following listing shows the mid-level IR when the inner loop over the input rows is unrolled by a factor of 2 and the two resulting tree walks are jammed together.


```

1 forest = ensemble(...)
2 for t = 0 to numTrees step 1 {
3   for i = 0 to batchSize step 2 {
4     tree = getTree(forest, t)
5     prediction1, prediction2 = InterleavedWalk((tree,
6       rows[i]), (tree, rows[i+1]))
7   }

```

When lowered to low-level IR, the operations to traverse each of the tree, input row pairs (the arguments to the `InterleavedWalk`) are interleaved. One step of the interleaved walk is listed below.

```

1 // ...
2 n1 = n2 = getRoot(tree)
3 // ...
4 threshold1 = loadThresholds(n1)
5 threshold2 = loadThresholds(n2)
6 featureIndex1 = loadFeatureIndices(n1)
7 featureIndex2 = loadFeatureIndices(n2)
8 feature1 = rows[i][featureIndex1]
9 feature2 = rows[i][featureIndex2]
10 pred1 = feature1 < threshold1
11 pred2 = feature2 < threshold2
12 n1 = getChildNode(n1, pred1)
13 n2 = getChildNode(n2, pred2)
14 // ...

```

These transformations are fairly general and are not aware of the in memory representation of the model. Therefore, they are reusable across different in memory representations - the ones that are currently built into Treebeard or ones that maybe added in the future. **TODO AP I feel the way this section is currently written makes the optimization seem extremely trivial. Is there a different way to present it?**

4.2 Loop peeling and walk unrolling

TREEBEARD splits the loop that performs a tree walk into two parts. As it is aware of the depth of the first leaf in a tree, it peels and introduces a prologue loop that walks down the tree a constant number of steps (up to first leaf) and then performs the rest of the tree walk in a separate loop. Further TREEBEARD unrolls the prologue completely, avoiding all traversal induced branching in it.

Note that for trees where TREEBEARD has already padded and balanced the tree (Section 3.5), walk unrolling completely avoids all traversal induced spills. **TODO kr: Add example?**

4.3 Parallelization

Currently, Treebeard performs a naive parallelization of the inference computation. When parallelism is enabled, the loop over the input rows is parallelized using OpenMP. Treebeard rewrites the mid-level IR by tiling the loop over the input rows with a tile size equal to the number of cores. As a concrete example, consider the case where we intend to perform inference using a model with 4 trees on a batch of 64 rows. Further, assume that we wish to parallelize this computation across 8 cores. Treebeard then generates the following IR.

```

1 forest = ensemble(...)
2 parallel.for i0 = 0 to 64 step 8 {
3   for i1 = 0 to 8 step 1 {
4     i = i0 + i1
5     prediction = 0
6     for t = 0 to 4 step 1 {
7       tree = getTree(forest, t)
8       treePrediction = WalkTree(tree, rows[i])
9       prediction = prediction + treePrediction
10    }

```

```

11 predictions[i] = prediction
12 }
13 }

```

Currently, as our focus is on single core performance, we do not do any parallelism related optimizations (such as loop tiling). We leave a more thorough exploration of parallelization to future work.

5. LOW LEVEL OPTIMIZATIONS

5.1 Vectorization

Vectorization performed by Treebeard is enabled by the tiling transformations described in section 3.2. When the low level IR is translated to LLVM IR, Treebeard generates LLVM instructions that operate on the thresholds and feature indices of nodes within a tile in a vector fashion. Therefore, thresholds and feature indices are loaded using vector loads and predicates are evaluated using vector comparisons. These vector LLVM IR instructions are then converted to vector instructions in the target ISA by the LLVM JIT.

The below listing shows some of the details of a vectorized tree walk.

```

1 // A lookup table that determines the child index of
2 // the next tile given the tile shape and the outcome
3 // of the vector comparison on the current tile
4 int16_t LUT[NUM_TILE_SHAPES, pow(2, TileSize)]
5
6 ResultType Prediction_Function(...) {
7   // ...
8   Node n = getRoot(tree)
9   while (isLeaf(tree, n) == false) do {
10     thresholds = loadThresholds(tree, n)
11     featureIndices = loadFeatureIndices(tree, n)
12     // Gather required feature from the current row
13     features = rows[i][featureIndices]
14     // Vector comparison of features and thresholds
15     comparison = features < thresholds
16
17     // Pack bits in comparison vector into an integer
18     comparisonIndex = combineBitsIntoInt(comparison)
19
20     // Get child index of tile we need to move to
21     tileShape = loadTileShape(tree, n)
22     childIndex = LUT[tileShapeID, comparisonIndex]
23
24     // Move to the correct child of the current node
25     n = getChildNode(tree, n, childIndex)
26   }
27   ThresholdType prediction = getLeafValue(n)
28   // ...
29 }

```

To evaluate the current tile, the vector of thresholds is first loaded (`loadThresholds`). This vector contains the thresholds of all nodes in the tile. Then, the features required for comparison are gathered into a vector (lines 11 and 13). The feature vector is compared to the threshold vector and the child tile to move to is determined (lines 15 to 25). More details about tile shapes and the look up table are presented in subsequent sections.

5.1.1 Tile Shapes

Informally, the *tile shape* is the shape of the region that encloses all nodes in a tile in a diagram of the decision tree. More formally, for a tile size n_t , each unique legal binary tree containing n_t nodes (nodes being indistinguishable) corresponds to a tile shape.

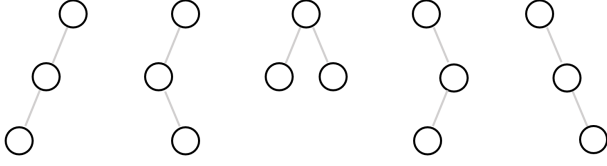


Figure 5: All possible tile shapes with tile size $n_t = 3$

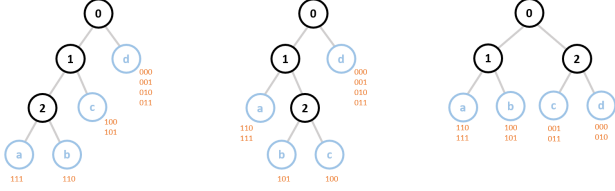


Figure 6: Example tile traversals with tile size $n_t = 3$

Figure 5 enumerates all tile shapes with a tile size of 3. There are a total of 5 tile shapes with size 3. The number of tile shapes with a tile size n_t , denoted by $NTS(n_t)$ is given by the following equation.

$$NTS(n) = \sum_{k=0}^{n-1} NTS(k) \times NTS(n-k-1) \quad (1)$$

where $NTS(0) = NTS(1) = 1$.

5.1.2 Tile Shapes and Decision Tree Inference

Treebeard uses vector instructions to accelerate decision tree walks. Vector instructions are used to evaluate the predicates of all the nodes in a tile simultaneously. However, once the predicates of all the nodes in the tile are evaluated, computing the next tile to move to, given the outcome of the comparison depends on the tile shape of the current tile. To illustrate this problem, consider the case of the tiles of size 3 shown in figure 6. The diagram shows 3 of the 5 possible tile shapes for a tile size of 3. The nodes drawn in black are members of the tile t_1 . The nodes in blue are the entry nodes of the children tiles of t_1 . **TODO Define entry nodes**

In the diagram, the bit strings (written in red) show which child we need to move to given the outcomes of the comparison. The bits represent the comparison outcomes of nodes and are in the order of the nodes in the tile – marked 0, 1 and 2 in the diagram, i.e., the MSB is the predicate outcome of node 0 and the LSB the predicate outcome of node 2. For example, for the first tile shape, if the predicates of all nodes are true (i.e. the comparison outcome is 111), the next node to evaluate is a . It is easy to see from the diagram that, depending on the tile shape, the same predicate outcomes can mean moving to different children. For example, for the outcome "011", the next tile is the 4th child (node d) for the first two tile shapes while it is the 3rd child for the other tile shape (node c).

5.1.3 Lookup Table

A lookup table (LUT) is used to solve the problem described in section 5.1.2, i.e. given the outcome of the comparisons of all nodes in a tile, determine the child tile we should

evaluate next. The LUT is indexed by the tile shape and the comparison outcome. Formally, the LUT is a map.

$$LUT : (TileShape, < Boolean \times n_t >) \rightarrow [0, n_t] \subset \mathbb{N}$$

where n_t is the tile size, $< Boolean \times n_t >$ is a vector of n_t booleans. The value returned by the LUT is the index of the child of the current tile that should be evaluated next. For example, if we are evaluating the first tile t in figure 6, and the result of the comparison is 110, then $LUT(TileShape(t), 110) = 1$ since the tile we need to evaluate next is the tile with node b , which is the second child of the current tile.

In order to realize this LUT in generated code, Treebeard associates a non-negative integer ID with every unique tile shape of the given tile size. The result of the comparison, a vector of booleans, can be interpreted as a 64-bit integer. Therefore, the LUT can be implemented as a 2 dimensional array. Treebeard computes the values in the LUT statically as the tile size is a compile time constant. **TODO AP What comes after subsection?**

5.2 In Memory Representation of Tiled Trees

Treebeard currently has two in memory representations for tiled trees - an array based representation and a sparse representation. Both representations use an array of structs to represent all tiles of the model.

5.2.1 Array Based Representation

Each tree in the model is represented as an array of tiles using the standard representation of trees as arrays. The root node is at index 0 and for a node at index n in the array, the index of its i^{th} child is given by $(n_t + 1)n + (i + 1)$ (nodes in the tree of tiles have $n_t + 1$ children). A tile is represented by an object of the following struct.

```

21 struct Tile {
22     // A vector of TileSize elements
23     <ThresholdType x TileSize> thresholds;
24     <FeatureIndexType x TileSize> featureIndices;
25     // Integer that identifies the tile shape
26     TileShapeIDType tileShapeID;
27 };

```

TODO AP Is this level of detail really needed? Also, the vector type notation needs to be introduced somewhere. Even though this representation is simple and efficient for small models, the memory required for bigger models is very large. This memory bloat causes performance problems because the span of the L1 TLB is not sufficient to efficiently translate addresses for the whole model. Storing leaves as full tiles (even though leaves just have to represent one value) and the empty space introduced due to the array based representation of trees that are not complete account for most of the increase.

5.2.2 Sparse Representation

The sparse representation tries to address the large memory footprint of the array based representation by doing the following.

- We add a child pointer to each tile to eliminate the wasted space in the array representation. This points to the first child of the tile. All children of a tile are stored contiguously.

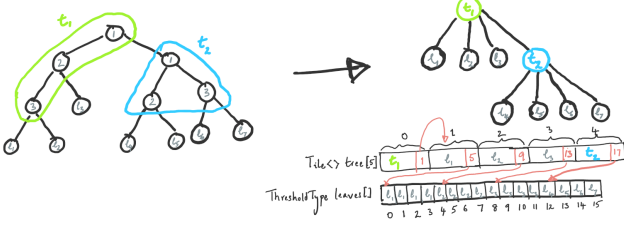


Figure 7: Sparse representation with tile size $n_t = 3$

- Leaves are stored as a separate array of scalar values. Across all our benchmarks, after tiling a large fraction of leaves are such that all their siblings are also leaves. Such leaves are directly moved into the leaves array. For leaves for which this property does not hold, an extra “hop” is added by making the original leaf tile a normal tile. All its children are made leaves with the same value as the original leaf.

Figure 7 shows some of the details of the sparse representation. The arrays depicted below show how the tree is represented in memory. The first array (`tree`) is an array of tiles and has 5 elements. Each element of the array represents a single tile and has the thresholds of the nodes, the feature indices, a tile shape ID and a pointer to the first child (shown explicitly in red).

As a specific example, consider the tile t_1 . The tile has four children – l_1, l_2, l_3 and t_2 in that order (left to right). These tiles are stored contiguously in the `tree` array and a pointer to the first of these, l_1 is stored in the tile t_1 (the index 1 is stored in the tile t_1 as shown).

Now consider the tile t_2 . Since all children of the tile t_2 are leaves, they are all moved into the `leaves` array. To store a pointer into the `leaves` array, we add $\text{len}(\text{tree})$ to the element index in the `leaves` array. The tile t_2 ’s child is the element at index 12 of the `leaves` array. Therefore, the index $12 + 5 = 17$ is stored in the tile t_2 . Any index i that is greater than the length of the `tree` array is regarded as an index into the `leaves` array. The index into the `leaves` array is $i - \text{len}(\text{tree})$.

The other aspect of the representation is that an extra hop is added for the leaves l_1, l_2 and l_3 in order to simplify code generation. This enforces the invariant that all leaves are stored in the `leaves` array and simplifies checking whether we’ve reached a leaf. Therefore, 4 new leaves are added as children for each of the original leaves l_1, l_2 and l_3 . Each of these 12 newly added leaves has the same value as its parent. These are the first 12 elements of the `leaves` array.

Even though we currently have implementations of the two representations detailed in sections 5.2.1 and 5.2.2, support for other representations is not hard to add. All optimizing passes that work on the high level and mid level IR will continue to work as is. Programmers need only provide new lowering passes for a few operations in the low level IR.

5.2.3 Code Generation for Probability Based Tiling

As probability based tiling pulls the most probable leaves of a decision tree nearest the root, it poses some implementation challenges. By design, the tiling process makes the tree of tiles imbalanced. The array based representation (section 5.2.1) cannot be used because of the memory footprint increase (a large part of the tree is empty, but would need to be allocated). On the other hand, the sparse representation in section 5.2.2 adds an extra hop for leaves that have non-leaf siblings. But this would mean that we add extra hops for the most probable leaves after probability based tiling which defeats the optimization. **TODO kr: connect to Peeling in MIR** We address these challenges using a code generation strategy. Treebeard peels the tree walk and specializes the leaf checks at higher levels to avoid the extra hop. Currently, we determine the maximum depth of leaves needed to cover 90 percent of the inputs and peel the tree walk by as many iterations. For example, consider the case where leaves until depth 2 are needed to cover 90 percent of the training input. Then, Treebeard generates the following IR.

```

28 // ...
29 tree = getTree(forest, t)
30 node = getRoot(tree)
31 node = traverseTreeTile(tree, node, rows[i])
32 if (isLeafTile(node)) {
33     treePrediction = getLeafValue(tree, node)
34 } else {
35     node = traverseTreeTile(tree, node, rows[i])
36     if (isLeafTile(node)) {
37         treePrediction = getLeafValue(tree, node)
38     } else {
39         // Loop based traversal
40     }
41 }
42 treePrediction = getLeafValue(tree, node)
43 // ...

```

The if statements check whether a node is a leaf tile and hence avoid the extra hop. While walk peeling is used to improve the performance of probability based tiling by specializing leaf tests, the transformation is by itself general and can be used in different contexts. For example, it could be used to elide leaf checks until a depth d is reached if we know all leaves are at a depth greater than d .

One other issue that the code generator needs to handle is that walks of different trees in the same ensemble may need to be peeled to different depths. A strategy similar to what is used for uniform tiling is used to handle this. Trees are reordered so that all trees with equal peeling depth are grouped together and the loops in the IR are fussed so that tree walks for these groups of trees can be specialized differently.

REFERENCES

- [1] C. Lattner and V. Adve, “Llvm: a compilation framework for lifelong program analysis and transformation,” in *International Symposium on Code Generation and Optimization, 2004. CGO 2004.*, 2004, pp. 75–86.